

Data Path : Z:\SVOASRV\HPCHEM1\BNA_P\DATA\BP100919\
 Data File : BP000598.D
 Acq On : 10 Oct 2019 03:06
 Operator : JU
 Sample : PB123661BL
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 SBLK61

Quant Time: Oct 10 03:50:05 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_P\METHODS\SOM-EPA-BP100919MA.M
 Quant Title : SVOA CALIBRATION
 QLast Update : Wed Oct 09 17:20:45 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	6.75	152	230313	20.00	ng/ul	0.00
18) Naphthalene-d8	8.03	136	887183	20.00	ng/ul	0.00
36) Acenaphthene-d10	9.80	164	517349	20.00	ng/ul	0.00
62) Phenanthrene-d10	11.29	188	970004	20.00	ng/ul	0.00
78) Chrysene-d12	13.96	240	841265	20.00	ng/ul	0.00
86) Perylene-d12	15.44	264	930803	20.00	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
3) 1,4-Dioxane-d8	2.36	96	40518	7.30	ng/uL	-0.01
5) Phenol-d5	6.39	99	662910	37.31	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.43	67	336673	37.91	ng/ul	0.00
9) 2-Chlorophenol-d4	6.52	132	558621	37.32	ng/ul	0.00
13) 4-Methylphenol-d8	7.13	113	511287	37.67	ng/ul	0.00
19) Nitrobenzene-d5	7.31	128	260603	38.64	ng/ul	0.00
22) 2-Nitrophenol-d4	7.64	143	286170	39.85	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	7.88	165	506105	36.82	ng/ul	0.00
29) 4-Chloroaniline-d4	8.10	131	716022	46.84	ng/ul	0.00
44) Dimethylphthalate-d6	9.49	166	1411301	38.97	ng/ul	0.00
47) Acenaphthylene-d8	9.65	160	1672110	37.53	ng/ul	0.00
52) 4-Nitrophenol-d4	9.90	143	215850	38.69	ng/ul	0.00
58) Fluorene-d10	10.32	176	1184829	38.24	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	10.39	200	136606	29.57	ng/ul	0.00
71) Anthracene-d10	11.35	188	1748782	38.92	ng/ul	0.00
79) Pyrene-d10	12.73	212	1822001	38.30	ng/ul	0.00
90) Benzo(a)pyrene-d12	15.35	264	1779509	38.72	ng/ul	0.00

Target Compounds	Ovalue
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(#) = qualifier out of range (m) = manual integration (+) = signals summed

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